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TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * Welcome to STN International * * * * * * * * *

NEWS 1 Web Page for STN Seminar Schedule - N. America
NEWS 2 APR 04 STN AnaVist, Version 1, to be discontinued
NEWS 3 APR 15 WPIDS, WPINDEX, and WPIX enhanced with new predefined hit display formats
NEWS 4 APR 28 EMBASE Controlled Term thesaurus enhanced
NEWS 5 APR 28 IMSRESEARCH reloaded with enhancements
NEWS 6 MAY 30 INPAFAMDB now available on STN for patent family searching
NEWS 7 MAY 30 DGENE, PCTGEN, and USGENE enhanced with new homology sequence search option
NEWS 8 JUN 06 EPFULL enhanced with 260,000 English abstracts
NEWS 9 JUN 06 KOREAPAT updated with 41,000 documents
NEWS 10 JUN 13 USPATFULL and USPAT2 updated with 11-character patent numbers for U.S. applications
NEWS 11 JUN 19 CAS REGISTRY includes selected substances from web-based collections
NEWS 12 JUN 25 CA/CAplus and USPAT databases updated with IPC reclassification data
NEWS 13 JUN 30 AEROSPACE enhanced with more than 1 million U.S. patent records
NEWS 14 JUN 30 EMBASE, EMBAL, and LEMBASE updated with additional options to display authors and affiliated organizations
NEWS 15 JUN 30 STN on the Web enhanced with new STN AnaVist Assistant and BLAST plug-in
NEWS 16 JUN 30 STN AnaVist enhanced with database content from EPFULL
NEWS 17 JUL 28 CA/CAplus patent coverage enhanced
NEWS 18 JUL 28 EPFULL enhanced with additional legal status information from the epoline Register
NEWS 19 JUL 28 IFICDB, IFIPAT, and IFIUDB reloaded with enhancements
NEWS 20 JUL 28 STN Viewer performance improved
NEWS 21 AUG 01 INPADOCDB and INPAFAMDB coverage enhanced
NEWS 22 AUG 13 CA/CAplus enhanced with printed Chemical Abstracts page images from 1967-1998
NEWS 23 AUG 15 CAOLD to be discontinued on December 31, 2008
NEWS 24 AUG 15 CAplus currency for Korean patents enhanced
NEWS 25 AUG 25 CA/CAplus, CASREACT, and IFI and USPAT databases enhanced for more flexible patent number searching
NEWS 26 AUG 27 CAS definition of basic patents expanded to ensure comprehensive access to substance and sequence information

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS STN Operating Hours Plus Help Desk Availability

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FILE 'HOME' ENTERED AT 11:42:36 ON 08 SEP 2008

=> file reg
COST IN U.S. DOLLARS
SINCE FILE ENTRY TOTAL
SESSION
FULL ESTIMATED COST 0.21 0.21

FILE 'REGISTRY' ENTERED AT 11:42:45 ON 08 SEP 2008
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STRUCTURE FILE UPDATES: 7 SEP 2008 HIGHEST RN 1047406-12-1
DICTIONARY FILE UPDATES: 7 SEP 2008 HIGHEST RN 1047406-12-1

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TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

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<http://www.cas.org/support/stn/gen/stndoc/properties.html>

=>
Uploading C:\Program Files\STNEXP\Queries\10596745s8.str



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chain nodes :
2 3 4 6 7 9 10 12 13 14 15 22
ring/chain nodes :
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chain bonds :
1-10 2-3 2-22 3-13 4-13 4-14 6-7 6-14 7-9 9-15 10-22 12-15
exact/norm bonds :
1-10 4-13 4-14 6-7 6-14 7-9 9-15 10-22 12-15
exact bonds :
2-3 2-22 3-13

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G1:C,O,N,Cy

G2:C,O,N,Cy

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Match level :
1:CLASS 2:CLASS 3:CLASS 4:Atom 6:CLASS 7:CLASS 9:CLASS 10:Atom 12:Atom
13:CLASS 14:CLASS 15:CLASS 22:CLASS
Element Count :
Node 10: Limited
    C,C4
    N,N2

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L1 STRUCTURE UPLOADED

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=> s l1 sss sam
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SAMPLE SCREEN SEARCH COMPLETED - 15463 TO ITERATE

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12.9% PROCESSED     2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

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2 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 301811 TO 316709
PROJECTED ANSWERS: 74 TO 544

L2 2 SEA SSS SAM L1

=> s c6n2/rf
L3 52023 C6N2/RF

=> s c4n2/rf
L4 2537685 C4N2/RF

=> s l1 sss sub=13 sam
SAMPLE SUBSET SEARCH INITIATED 11:44:46 FILE 'REGISTRY'
SAMPLE SUBSET SCREEN SEARCH COMPLETED - 19 TO ITERATE

100.0% PROCESSED 19 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **COMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 119 TO 641
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 0 TO 0

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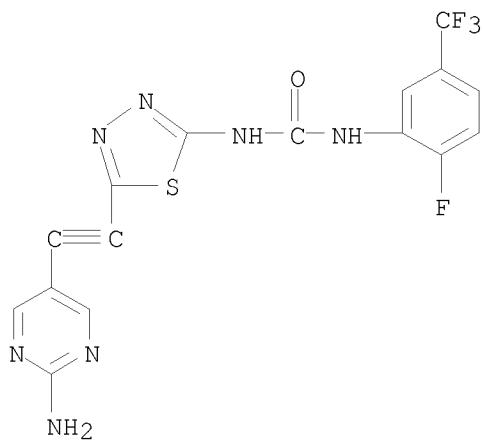
100.0% PROCESSED 1287 ITERATIONS 15 ANSWERS
SEARCH TIME: 00.00.01

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PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 68 TO 532

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=> d scan

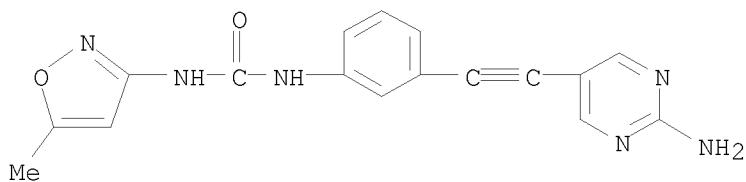
L6 15 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-1,3,4-thiadiazol-2-yl]-N'-[2-fluoro-5-(trifluoromethyl)phenyl]-
MF C16 H9 F4 N7 O S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

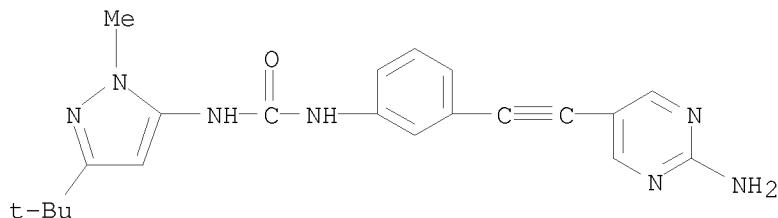
L6 15 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-methyl-3-isoxazolyl)-
 MF C17 H14 N6 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L6 15 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-
 MF C21 H23 N7 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s 11 sss sub=14 full
FULL SUBSET SEARCH INITIATED 11:46:00 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 26775 TO ITERATE

100.0% PROCESSED 26775 ITERATIONS 317 ANSWERS
SEARCH TIME: 00.00.01

L7 317 SEA SUB=L4 SSS FUL L1

=> s 11 sss full
FULL SEARCH INITIATED 11:46:22 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 309531 TO ITERATE

100.0% PROCESSED 309531 ITERATIONS 317 ANSWERS
SEARCH TIME: 00.00.03

L8 317 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL
 ENTRY SESSION
FULL ESTIMATED COST 368.86 369.07

FILE 'CAPLUS' ENTERED AT 11:46:37 ON 08 SEP 2008
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FILE COVERS 1907 - 8 Sep 2008 VOL 149 ISS 11
FILE LAST UPDATED: 7 Sep 2008 (20080907/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

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=> s 18 and 17
 9 L8
 9 L7
L9 9 L8 AND L7

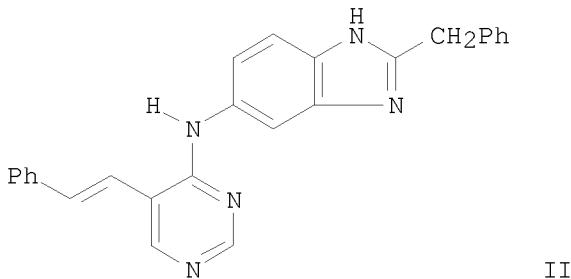
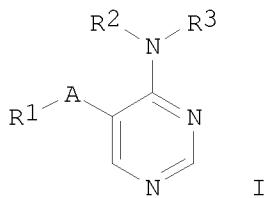
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L9 ANSWER 9 OF 9 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:158661 CAPLUS
 DOCUMENT NUMBER: 142:240460
 TITLE: Preparation of pyrimidine derivatives as ErbB kinase inhibitors
 INVENTOR(S): Reno, Michael John; Stevens, Kirk Lawrence; Waterson, Alex Gregory; Zhang, Yuemei
 PATENT ASSIGNEE(S): Smithkline Beecham Corporation, USA
 SOURCE: PCT Int. Appl., 132 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005016914	A1	20050224	WO 2004-US26251	20040811
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP 1654251	A1	20060510	EP 2004-781004	20040811
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, HR				
JP 2007502298	T	20070208	JP 2006-523388	20040811
US 20060205740	A1	20060914	US 2006-568052	20060210
PRIORITY APPLN. INFO.:			US 2003-495180P	P 20030814
			WO 2004-US26251	W 20040811

OTHER SOURCE(S): CASREACT 142:240460; MARPAT 142:240460
 GI



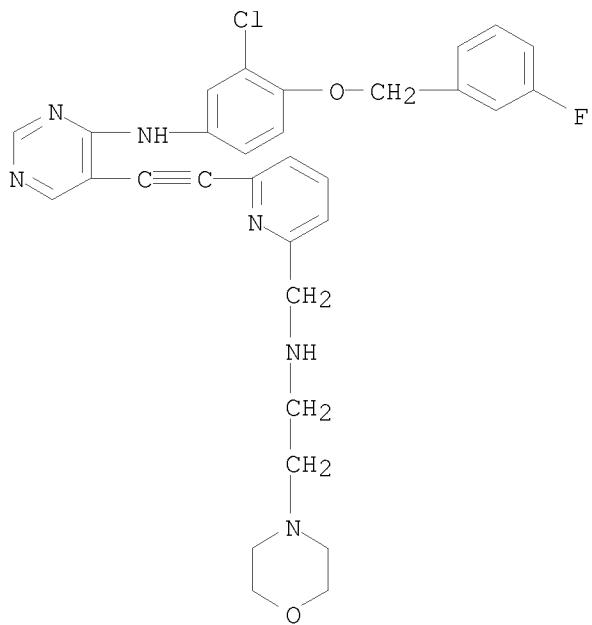
AB Title compds. I [wherein A = alkenylene, alkynylene; R = alkylene; R₁ = -(Z)-(Z₁)_m-(Z₂)_n; Z = hetero/aryl, hetero/arylene; Z₁ = CH₂ where m = 0-1; Z₂ = OH and derivs., halo, CN, CONH₂ and derivs. or heterocyclyl, where n = 0-1, etc.; R₂ = H, alkyl; R₃ = -(Q)-(Q₁)_r-(Q₂); Q = hetero/arylene; Q₁ = O, where r = 0-1; Q₂ = arylalkyl, hetero/aryl; and their salts, solvates, and physiol. functional derivs.] were prepared as ErbB kinase inhibitors for treating cancer. Thus, reacting 2-benzyl-N-(5-vinylpyrimidin-4-yl)-1H-benzimidazol-5-amine (preparation given) with Ph iodide gave pyrimidine II in 8%. I showed inhibitory activity vs. EGFR, ErbB-2, and ErbB-4 protein tyrosine kinases with a pIC₅₀ ≥ 5.0. I are useful in the treatment of diseases associated with inappropriate ErbB family kinase activity.

IT 845657-86-5P, N-[3-Chloro-4-[(3-fluorobenzyl)oxy]phenyl]-5-[[6-[[[2-(4-morpholinyl)ethyl]amino]methyl]-2-pyridinyl]ethynyl]-4-pyrimidinamine 845657-90-1P, N-[3-Chloro-4-[(3-fluorobenzyl)oxy]phenyl]-5-[[6-[[[3-(1H-imidazol-1-yl)propyl]amino]methyl]-2-pyridinyl]ethynyl]-4-pyrimidinamine 845658-10-8P,
1-[[6-[[4-[3-Chloro-4-[(3-fluorobenzyl)oxy]anilino]-5-pyrimidinyl]ethynyl]-2-pyridinyl]methyl]-3-[2-(4-morpholinyl)ethyl]urea
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug candidate; preparation of pyrimidines as ErB kinase inhibitors)

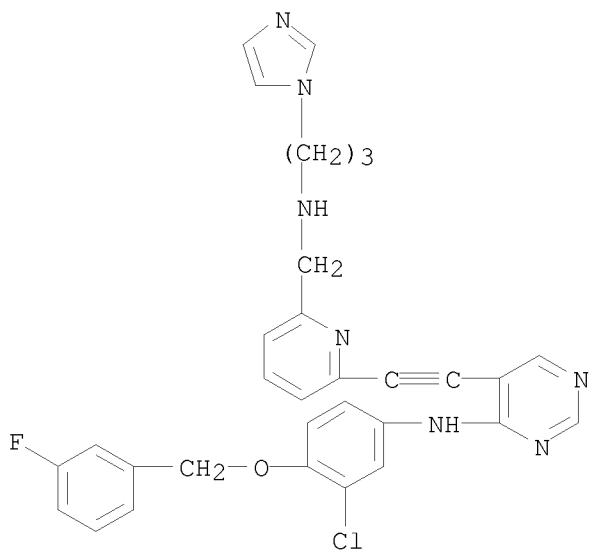
RN 845657-86-5 CAPLUS

CN 4-Morpholineethanamine, N-[[6-[2-[4-[(3-chloro-4-[(3-fluorophenyl)methoxy]phenyl)amino]-5-pyrimidinyl]ethynyl]-2-pyridinyl]methyl]- (CA INDEX NAME)



RN 845657-90-1 CAPLUS

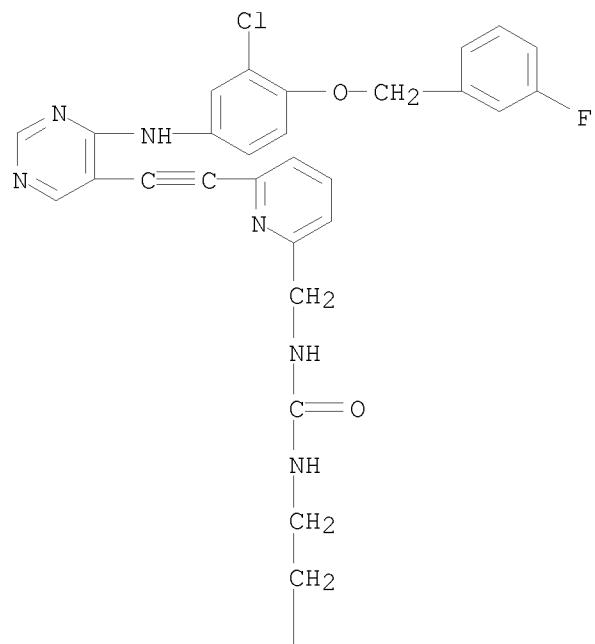
CN 4-Pyrimidinamine, N-[3-chloro-4-[(3-fluorophenyl)methoxy]phenyl]-5-[2-[6-[[3-(1H-imidazol-1-yl)propyl]amino]methyl]-2-pyridinyl]ethynyl]- (CA INDEX NAME)



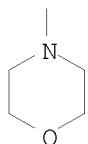
RN 845658-10-8 CAPLUS

CN Urea, N-[6-[2-[4-[3-chloro-4-[(3-fluorophenyl)methoxy]phenyl]amino]-5-pyrimidinyl]ethynyl]-2-pyridinyl]methyl]-N'-(2-(4-morpholinyl)ethyl)- (CA INDEX NAME)

PAGE 1-A



PAGE 2-A



REFERENCE COUNT: 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

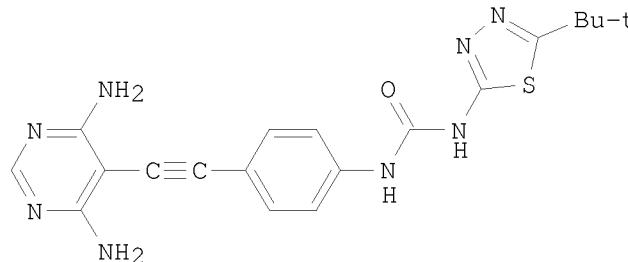
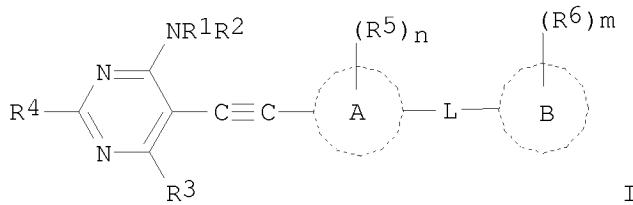
=> d ibib abs hitstr 8

L9 ANSWER 8 OF 9 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2005:588667 CAPLUS
DOCUMENT NUMBER: 143:115556
TITLE: Preparation of 4-aminopyrimidine derivatives as inhibitors of Tie2 receptor tyrosine kinases
INVENTOR(S): Jones, Clifford David; Luke, Richard William Arthur; McCoull, William
PATENT ASSIGNEE(S): Astrazeneca AB, Swed.; Astrazeneca UK Limited
SOURCE: PCT Int. Appl., 129 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005060969	A1	20050707	WO 2004-GB5332	20041220

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH,
 CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
 GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,
 LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
 NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
 TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
 RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM,
 AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,
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 RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML,
 MR, NE, SN, TD, TG
 EP 1737462 A1 20070103 EP 2004-806134 20041220
 EP 1737462 B1 20080730
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 IS, IT, LI, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR
 CN 1917880 A 20070221 CN 2004-80041936 20041220
 JP 2007517006 T 20070628 JP 2006-546305 20041220
 US 20080027076 A1 20080131 US 2006-596740 20060622
 IN 2006MN00847 A 20070420 IN 2006-MN847 20060717
 PRIORITY APPLN. INFO.: GB 2003-30001 A 20031224
 GB 2004-16850 A 20040729
 WO 2004-GB5332 W 20041220

OTHER SOURCE(S): CASREACT 143:115556; MARPAT 143:115556
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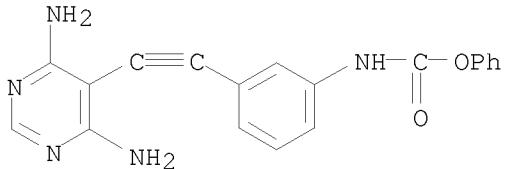


AB Title compds. I [wherein R1, R2 = H, alkyl, alkanoyl; R3, R4 = H, alkyl, alkoxy; R5 = cyclopropyl, halo, cyano; m, n = 0-3; R6 = halo, oxo, cyano; etc., or salts thereof] were prepared as inhibitors of Tie2 receptor tyrosine kinases. Processes for the synthesis of I and some intermediates involved are claimed. For example, urea II was synthesized in 21% yield by condensation of the corresponding aniline with Ph thiadiazolylcarbamate in the presence of Et3N in THF under microwave irradiation. This urea showed inhibition against Tie2 receptor tyrosine kinase in vitro and inhibition of autophosphorylation of Tie2 receptor tyrosine kinase with IC50 values of 0.879 μM and 5.557 μM, resp. Therefore, I and their pharmaceutical compns. have potential use in the production of an anti-angiogenic effect in a warm-blooded animal.
 IT 857287-13-9P, Phenyl [3-[(4,6-diaminopyrimidin-5-

yl)ethynyl]phenyl]carbamate 857287-53-7P, N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-(3-[4-[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl)ethynyl]phenyl]urea
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
 (inhibitor; preparation of aminopyrimidine derivs. as inhibitors of Tie2 receptor tyrosine kinases)

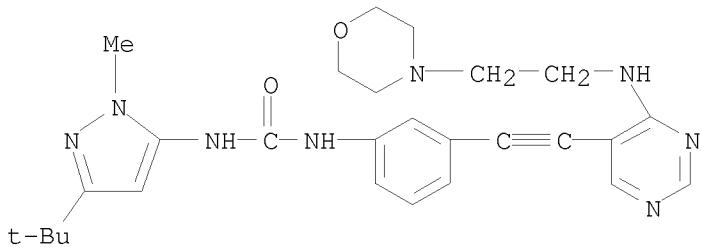
RN 857287-13-9 CAPLUS

CN Carbamic acid, [3-[(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



RN 857287-53-7 CAPLUS

CN Urea, N-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-N'-(3-[2-[4-[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl)- (CA INDEX NAME)



IT 857287-02-6P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-phenylurea 857287-04-8P, 2-Phenyl-N-[3-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl]acetamide 857287-05-9P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3,4-dichlorophenyl)urea 857287-06-0P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2-(trifluoromethyl)phenyl)urea 857287-07-1P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-(trifluoromethyl)phenyl)urea 857287-08-2P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)phenyl)urea 857287-09-3P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2-fluoro-5-(trifluoromethyl)phenyl)urea 857287-10-6P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methoxyphenyl)urea 857287-11-7P, Phenyl [4-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl]carbamate 857287-14-0P, N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-(4-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl)urea 857287-15-1P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisothiazol-5-yl)urea 857287-16-2P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisoxazol-5-yl)urea 857287-17-3P, N-[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)pyridin-2-yl)urea 857287-18-4P, N-[3-[[4-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]amino]carbonyl]amino]phenyl]acetamide 857287-19-5P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisothiazol-5-yl)urea 857287-20-8P, N-[3-[[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]amino]carbonyl]ami

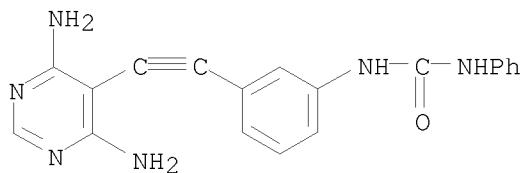
no]phenyl]acetamide 857287-21-9P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)pyridin-2-yl)urea
857287-22-0P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(3-methylisoxazol-5-yl)urea 857287-23-1P, N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-(3-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl)urea
857287-24-2P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4,6-diaminopyrimidin-5-yl)ethynyl]phenyl)urea 857287-25-3P,
N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2,3-dihydro-1,4-benzodioxin-6-yl)urea 857287-26-4P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2-(morpholin-4-yl)phenyl)urea
857287-27-5P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(1-methylpiperidin-4-yl)urea 857287-28-6P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(1-propylpiperidin-4-yl)urea
857287-29-7P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-2-(2-methoxyphenyl)acetamide 857287-30-0P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-2-[3-(trifluoromethyl)phenyl]acetamide 857287-31-1P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-2-[4-(trifluoromethyl)phenyl]acetamide
857287-32-2P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-2-(3-methoxyphenyl)acetamide 857287-35-5P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-(methylamino)pyrimidin-5-yl)ethynyl]phenyl)urea
857287-36-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[3-(isopropylamino)propyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-37-7P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[2-(pyrrolidin-1-yl)ethyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-38-8P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[5-tert-butylisoxazol-3-yl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-39-9P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[3-(dimethylamino)propyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-40-2P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[2-hydroxyethyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea 857287-41-3P
, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[2-(morpholin-4-yl)ethyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea 857287-42-4P,
N-[3-[(4-[(4-Aminobutyl)amino)pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)-N'-(3-[(4-[(2-
y1)-N'-[3-[(4-[(3-(pyrrolidin-1-yl)propyl]amino)pyrimidin-5-yl]ethynyl]phenyl)urea 857287-44-6P, N-(5-tert-Butylisoxazol-3-
y1)-N'-(3-[(4-[(2,4-dimethoxybenzyl)amino)pyrimidin-5-yl]ethynyl]phenyl)urea 857287-45-7P, N-[3-[(4-[(2-
Aminoethyl)amino)pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)urea 857287-46-8P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-
[[2-(dimethylamino)ethyl]amino)pyrimidin-5-yl]ethynyl]phenyl)urea
857287-47-9P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[
(dimethylamino)butyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-48-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[N-[2-
(dimethylamino)ethyl]methylamino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-49-1P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[
(piperidin-1-yl)ethyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-50-4P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[
(morpholin-4-yl)propyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-51-5P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[
(piperidin-1-yl)propyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-52-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-(3-[(4-[
methylpiperazin-1-yl)propyl]amino)pyrimidin-5-yl]ethynyl)phenyl)urea
857287-57-1P, N-[3-[(4,6-Diaminopyrimidin-5-yl)ethynyl]phenyl]-N'-(2,3-dihydro-1H-inden-1-yl)urea

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(inhibitor; preparation of aminopyrimidine derivs. as inhibitors of Tie2 receptor tyrosine kinases)

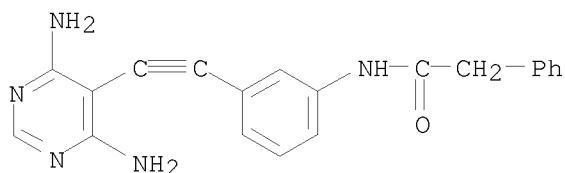
RN 857287-02-6 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-phenyl- (CA INDEX NAME)



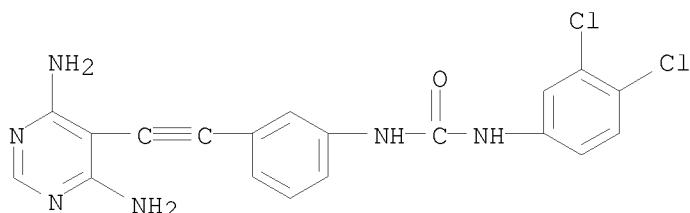
RN 857287-04-8 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]- (CA INDEX NAME)



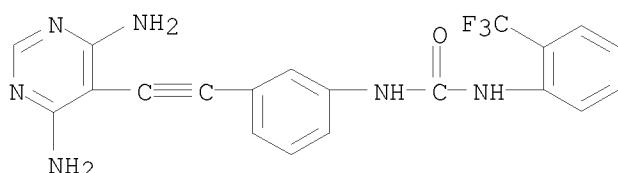
RN 857287-05-9 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3,4-dichlorophenyl)- (CA INDEX NAME)



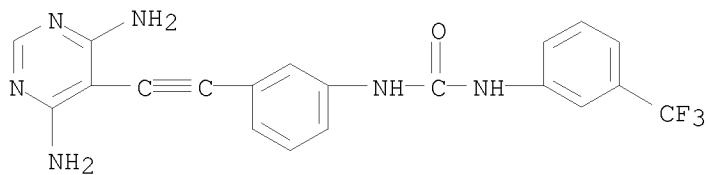
RN 857287-06-0 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2-(trifluoromethyl)phenyl)- (CA INDEX NAME)

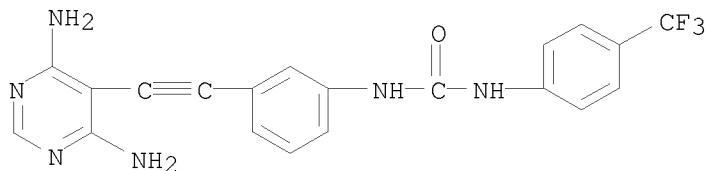


RN 857287-07-1 CAPLUS

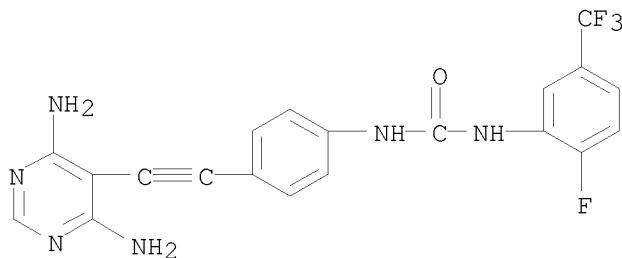
CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-(trifluoromethyl)phenyl)- (CA INDEX NAME)



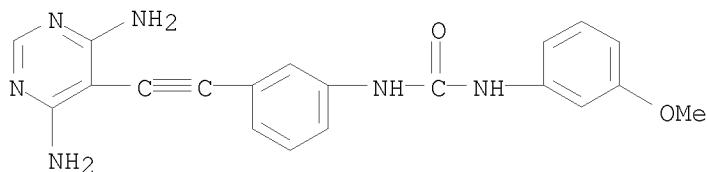
RN 857287-08-2 CAPLUS
 CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)phenyl)- (CA INDEX NAME)



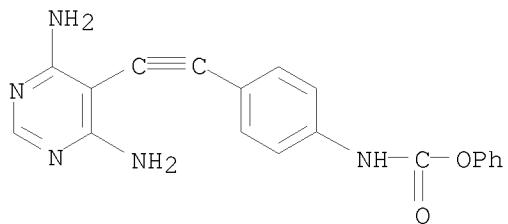
RN 857287-09-3 CAPLUS
 CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2-fluoro-5-(trifluoromethyl)phenyl)- (CA INDEX NAME)



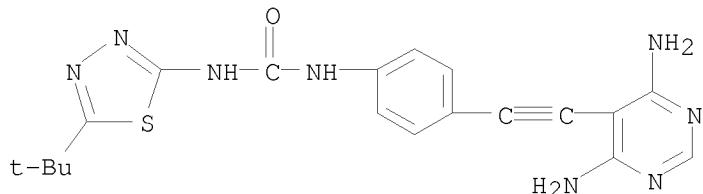
RN 857287-10-6 CAPLUS
 CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methoxyphenyl)- (CA INDEX NAME)



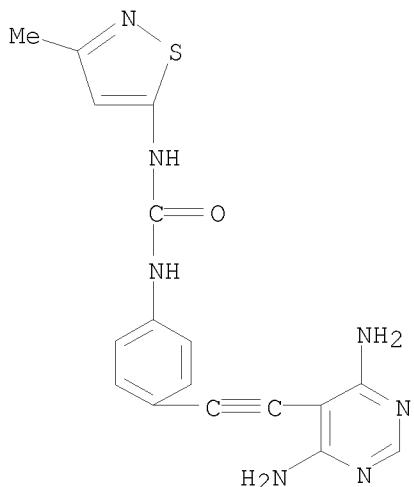
RN 857287-11-7 CAPLUS
 CN Carbamic acid, [4-[(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



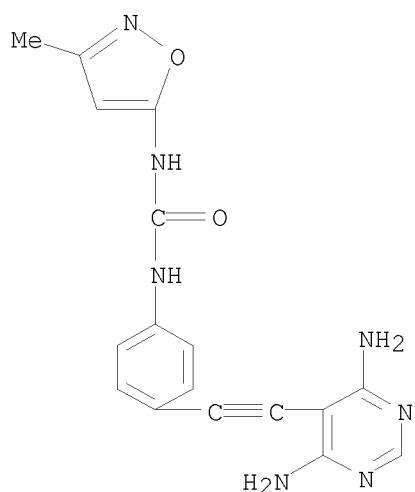
RN 857287-14-0 CAPLUS
 CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl)- (CA INDEX NAME)



RN 857287-15-1 CAPLUS
 CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isothiazolyl)- (CA INDEX NAME)

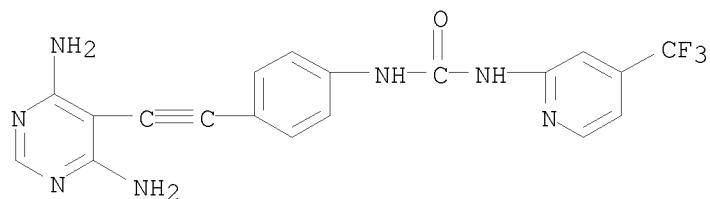


RN 857287-16-2 CAPLUS
 CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isoxazolyl)- (CA INDEX NAME)



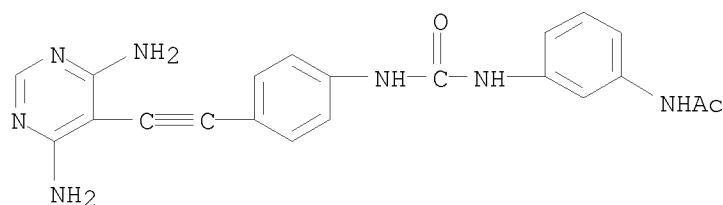
RN 857287-17-3 CAPLUS

CN Urea, N-[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)-2-pyridinyl)- (CA INDEX NAME)



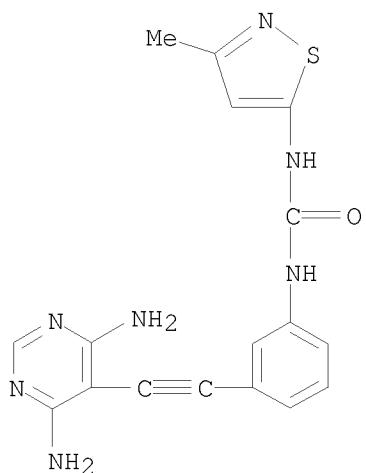
RN 857287-18-4 CAPLUS

CN Acetamide, N-[3-[[[4-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]amino]carbonyl]amino]phenyl- (CA INDEX NAME)



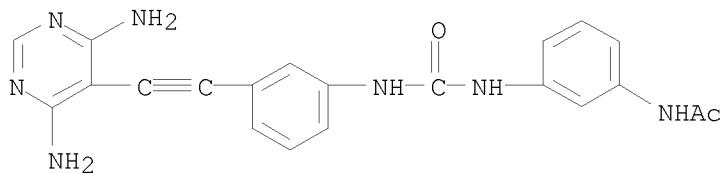
RN 857287-19-5 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isothiazolyl)- (CA INDEX NAME)



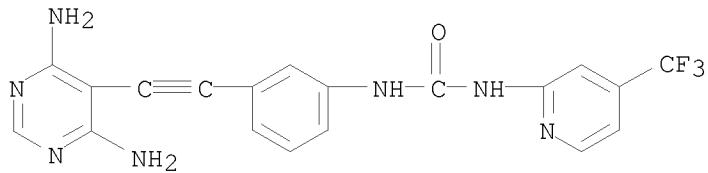
RN 857287-20-8 CAPLUS

CN Acetamide, N-[3-[[[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]amino]carbonyl]amino]phenyl]- (CA INDEX NAME)



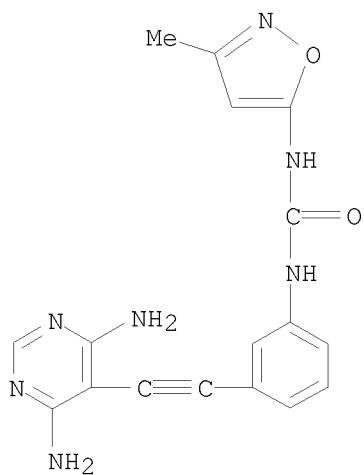
RN 857287-21-9 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(4-(trifluoromethyl)-2-pyridinyl)- (CA INDEX NAME)



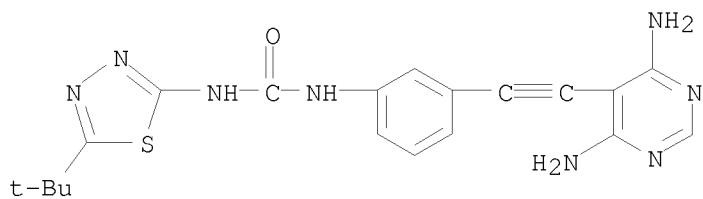
RN 857287-22-0 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-methyl-5-isoxazolyl)- (CA INDEX NAME)



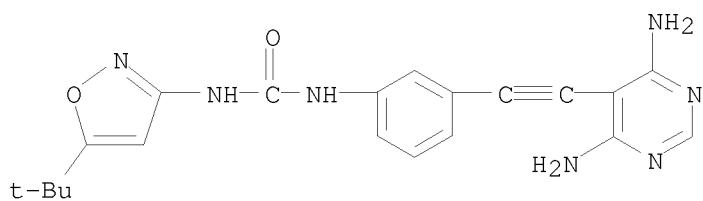
RN 857287-23-1 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl)- (CA INDEX NAME)



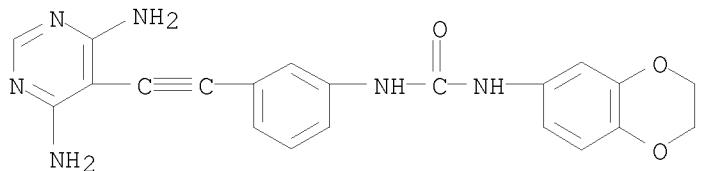
RN 857287-24-2 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)



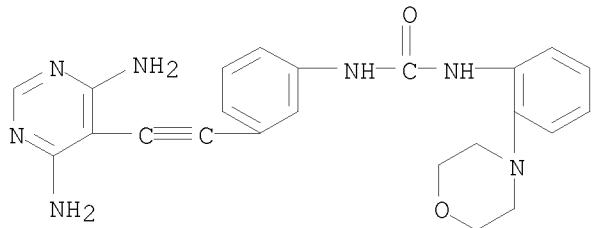
RN 857287-25-3 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2,3-dihydro-1,4-benzodioxin-6-yl)- (CA INDEX NAME)



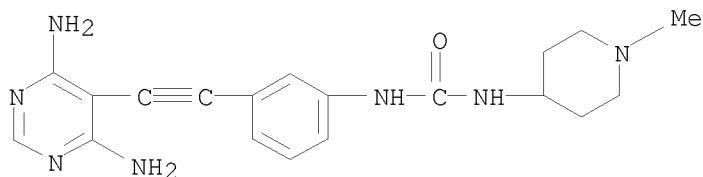
RN 857287-26-4 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2-(4-morpholinyl)phenyl)- (CA INDEX NAME)



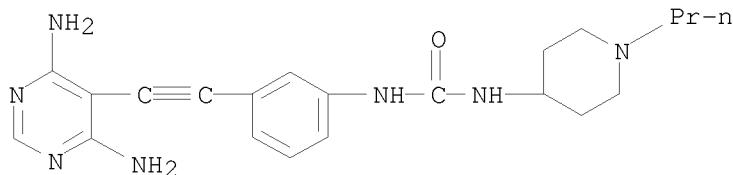
RN 857287-27-5 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-methyl-4-piperidinyl)- (CA INDEX NAME)



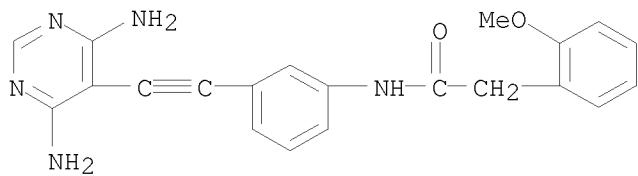
RN 857287-28-6 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-propyl-4-piperidinyl)- (CA INDEX NAME)



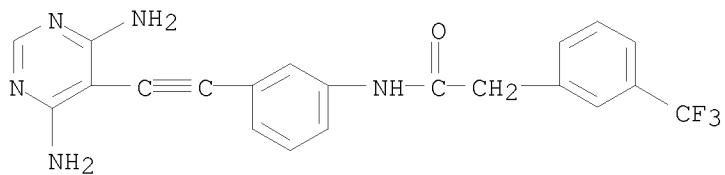
RN 857287-29-7 CAPLUS

CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-2-methoxy- (CA INDEX NAME)

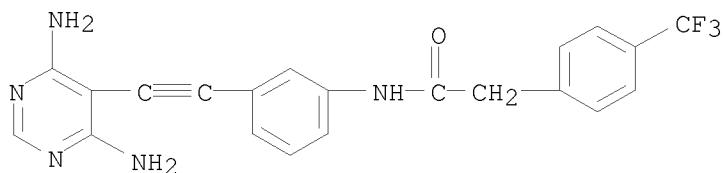


RN 857287-30-0 CAPLUS

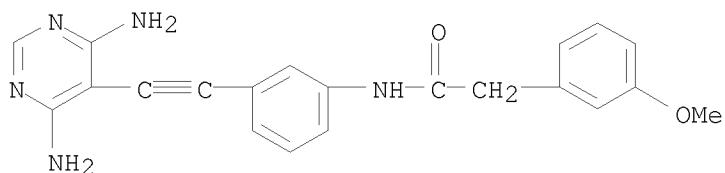
CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-3-(trifluoromethyl)- (CA INDEX NAME)



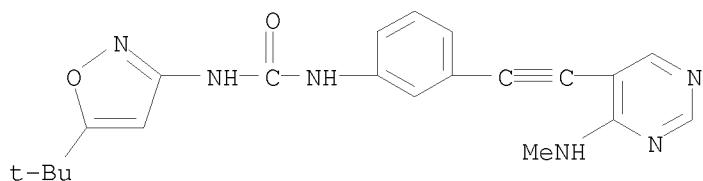
RN 857287-31-1 CAPLUS
 CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-4-(trifluoromethyl)- (CA INDEX NAME)



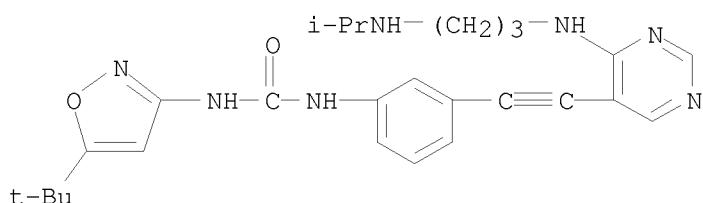
RN 857287-32-2 CAPLUS
 CN Benzeneacetamide, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-3-methoxy- (CA INDEX NAME)



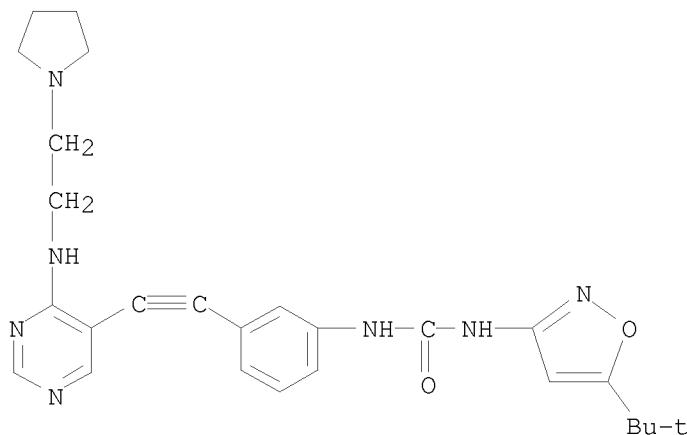
RN 857287-35-5 CAPLUS
 CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-(methylamino)-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



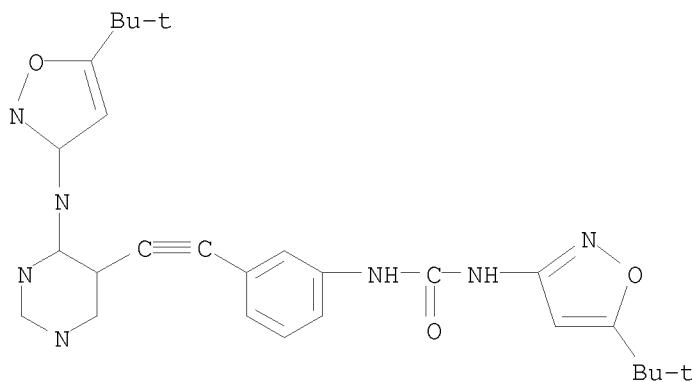
RN 857287-36-6 CAPLUS
 CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[3-[(1-methylethyl)amino]propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



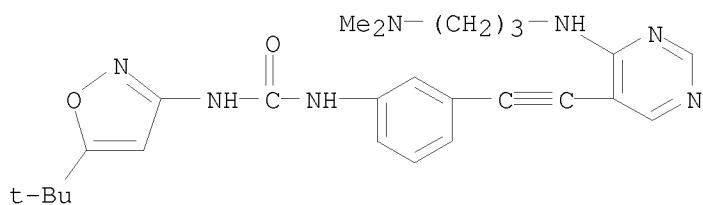
RN 857287-37-7 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857287-38-8 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

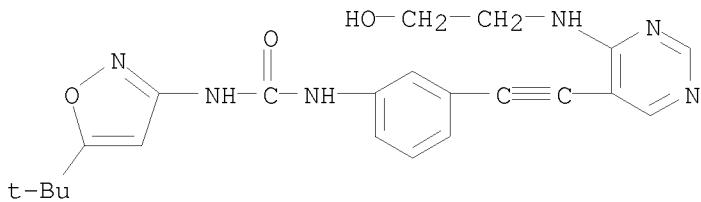


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
RN 857287-39-9 CAPLUS
CN Urea, N-[3-[2-[4-[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



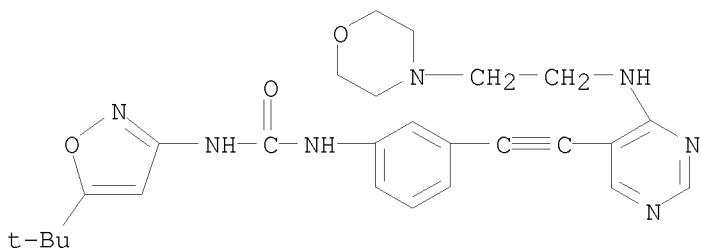
RN 857287-40-2 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[2-

hydroxyethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



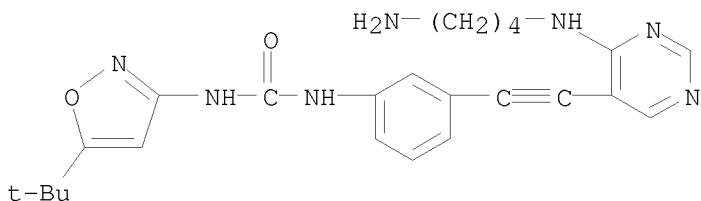
RN 857287-41-3 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[2-[4-morpholinyl]ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



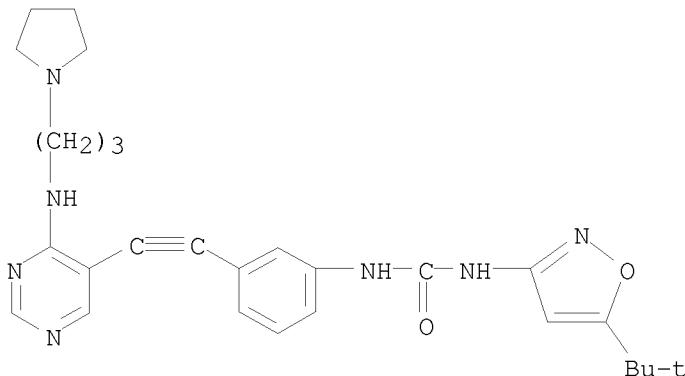
RN 857287-42-4 CAPLUS

CN Urea, N-[3-[2-[4-[(4-aminobutyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



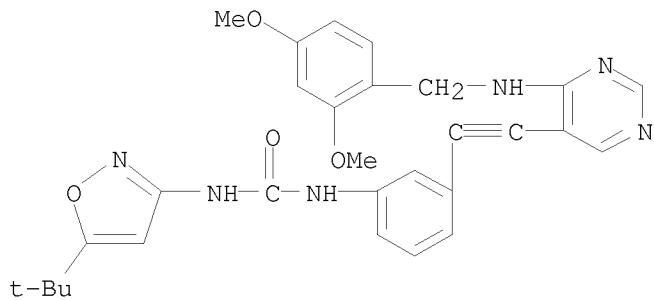
RN 857287-43-5 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[3-(1-pyrrolidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



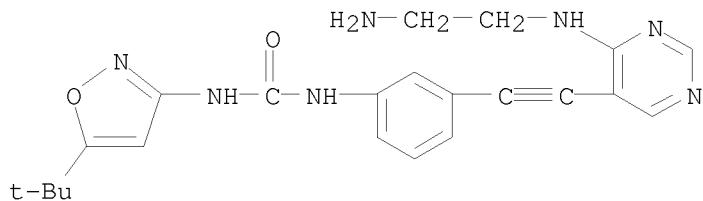
RN 857287-44-6 CAPLUS

CN Urea, N-[3-[2-[4-[(2,4-dimethoxyphenyl)methyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)



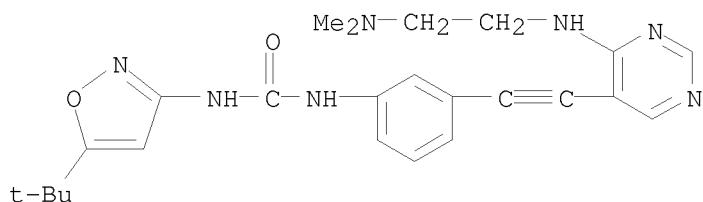
RN 857287-45-7 CAPLUS

CN Urea, N-[3-[2-[4-[(2-aminoethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)



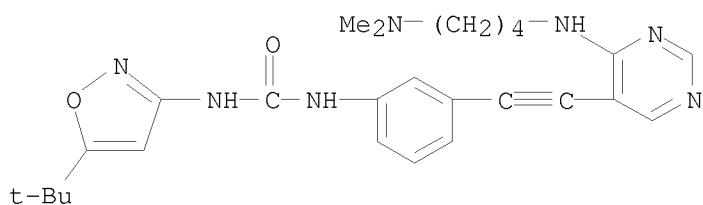
RN 857287-46-8 CAPLUS

CN Urea, N-[3-[2-[4-[(2-(dimethylamino)ethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)

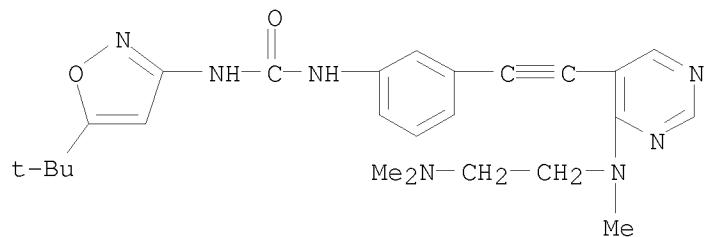


RN 857287-47-9 CAPLUS

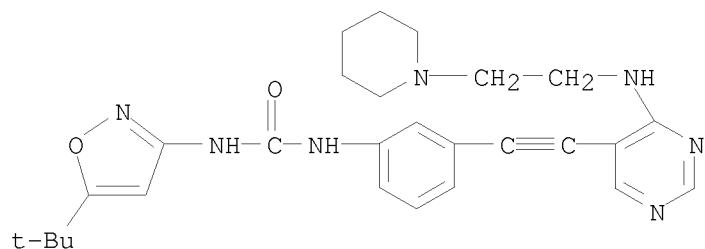
CN Urea, N-[3-[2-[4-[(4-(dimethylamino)butyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)



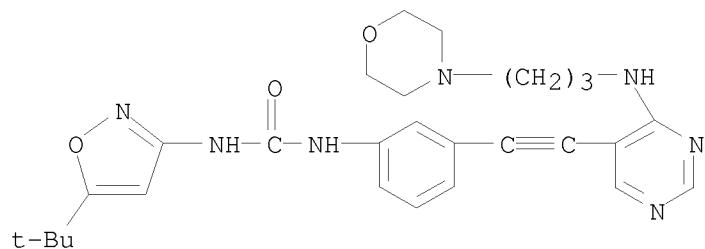
RN 857287-48-0 CAPLUS
CN Urea, N-[3-[2-[4-[[2-(dimethylamino)ethyl]methylamino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-(1,1-dimethylethyl)-3-isoxazolyl)- (CA INDEX NAME)



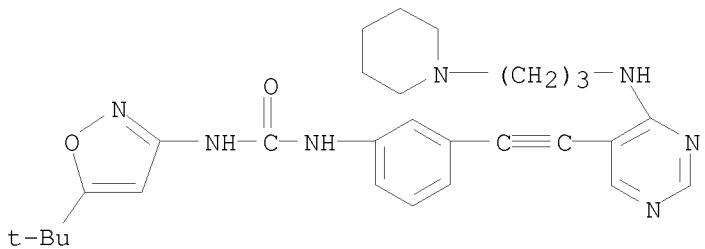
RN 857287-49-1 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-(3-[2-[4-[[2-(1-piperidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl)- (CA INDEX NAME)



RN 857287-50-4 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-(3-[2-[4-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl)- (CA INDEX NAME)

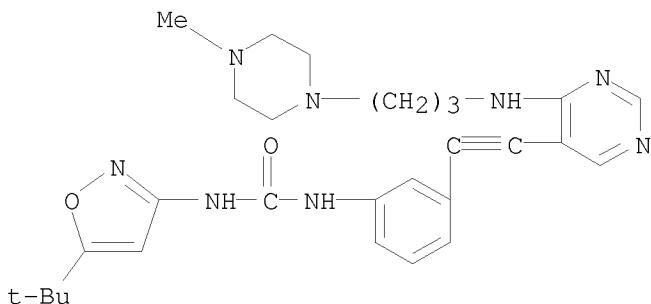


RN 857287-51-5 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-(3-[2-[4-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl)- (CA INDEX NAME)



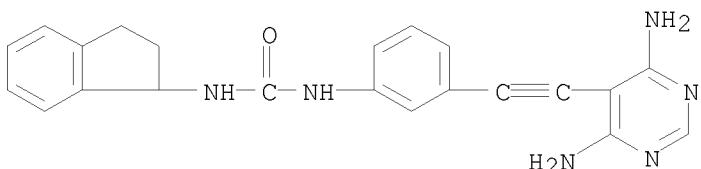
RN 857287-52-6 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[4-[3-(4-methyl-1-piperazinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857287-57-1 CAPLUS

CN Urea, N-[3-[2-(4,6-diamino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2,3-dihydro-1H-inden-1-yl)- (CA INDEX NAME)



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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SINCE FILE	TOTAL
ENTRY	SESSION

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